STN Columbus

* * *	* *	* *	* *	* Welcome to STN International * * * * * * * * *
NEWS NEWS	1 2	APR	02	Web Page for STN Seminar Schedule - N. America CAS Registry Number Crossover Limits Increased to 500,000 in Key STN Databases
NEWS	3	APR	02	PATDPAFULL: Application and priority number formats enhanced
NEWS	4	APR		DWPI: New display format ALLSTR available
NEWS		APR		New Thesaurus Added to Derwent Databases for Smooth Sailing through U.S. Patent Codes
NEWS	6	APR		EMBASE Adds Unique Records from MEDLINE, Expanding Coverage back to 1948
NEWS	7	APR	07	CA/CAplus CLASS Display Streamlined with Removal of Pre-IPC 8 Data Fields
NEWS	8	APR	07	50,000 World Traditional Medicine (WTM) Patents Now Available in CAplus
NEWS		APR		MEDLINE Coverage Is Extended Back to 1947
NEWS	10	JUN	16	WPI First View (File WPIFV) will no longer be available after July 30, 2010
NEWS		JUN		DWPI: New coverage - French Granted Patents
NEWS	12	JUN	18	CAS and FIZ Karlsruhe announce plans for a new STN platform
NEWS	13	JUN	18	IPC codes have been added to the INSPEC backfile (1969-2009)
NEWS	14	JUN	21	Removal of Pre-IPC 8 data fields streamline displays in CA/CAplus, CASREACT, and MARPAT
NEWS	15	JUN	21	Access an additional 1.8 million records exclusively
				enhanced with 1.9 million CAS Registry Numbers EMBASE Classic on STN
NEWS	16	JUN	28	Introducing "CAS Chemistry Research Report": 40 Years
				of Biofuel Research Reveal China Now Atop U.S. in Patenting and Commercialization of Bioethanol
NEWS	17	JUN	29	Enhanced Batch Search Options in DGENE, USGENE, and PCTGEN
NEWS	18	JUL	19	Enhancement of citation information in INPADOC databases provides new, more efficient competitor analyses
NEWS EXPRESS FEBRUARY 15 10 CURRENT WINDOWS VERSION IS V8.4.2, AND CURRENT DISCOVER FILE IS DATED 15 JANUARY 2010.				
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FILE 'HOME' ENTERED AT 22:49:21 ON 20 JUL 2010				
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FILE 'REGISTRY' ENTERED AT 22:49:35 ON 20 JUL 2010				
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 19 JUL 2010 HIGHEST RN 1233120-12-1 DICTIONARY FILE UPDATES: 19 JUL 2010 HIGHEST RN 1233120-12-1

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TSCA INFORMATION NOW CURRENT THROUGH January 8, 2010.

Please note that search-term pricing does apply when conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

http://www.cas.org/support/stngen/stndoc/properties.html

```
=> e ezetimibe/cn
E1
                     EZENTIA/CN
              1
E2
                    EZESLIDE/CN
              1 --> EZETIMIBE/CN
E3
                    EZETIMIBE GLUCURONIDE/CN
E4
              1
E5
              1
                    EZETIMIBE MONOHYDRATE/CN
              1
                    EZETROL/CN
E.6
                    EZFENI/CN
E7
              1
E8
              1
                    EZG-60G/CN
                    EZH 2/CN
E9
              1
E10
              1
                    EZH1/CN
                    EZH2 PROTEIN (XENOPUS LAEVIS CLONE MGC: 79865 IMAGE: 5073002 G
E11
              1
                     ENE EZH2)/CN
E12
              1
                     EZH4/CN
=> s e3
              1 EZETIMIBE/CN
L1
=> d
1.1
     ANSWER 1 OF 1 REGISTRY COPYRIGHT 2010 ACS on STN
     163222-33-1 REGISTRY
     Entered STN: 24 May 1995
ED
     2-Azetidinone, 1-(4-fluorophenyl)-3-[(3S)-3-(4-fluorophenyl)-3-
CN
     hydroxypropyl]-4-(4-hydroxyphenyl)-, (3R, 4S)- (CA INDEX NAME)
OTHER CA INDEX NAMES:
     2-Azetidinone, 1-(4-fluorophenyl)-3-[3-(4-fluorophenyl)-3-hydroxypropyl]-4-
     (4-hydroxyphenyl)-, [3R-[3\alpha(S^*), 4\beta]]-
OTHER NAMES:
CN
     (-) -Sch 58235
CN
     1-(4-\text{Fluoropheny1})-3(R)-[3-(4-\text{fluoropheny1})-3(S)-\text{hydroxypropy1}]-4(S)-(4-\text{fluoropheny1})
     hydroxyphenyl)azetidin-2-one
CN
     Ezedoc
CN
     Ezentia
     Ezetimibe
CN
CN
     Ezetrol
CN
     Ezta
     Sch 58235
CN
CN
     Zetia
FS
     STEREOSEARCH
MF
     C24 H21 F2 N O3
CI
     COM
SR
LC
     STN Files:
                    ADISINSIGHT, AGRICOLA, ANABSTR, BIOSIS, BIOTECHNO, CA,
       CAPLUS, CASREACT, CBNB, CHEMCATS, CIN, EMBASE, HSDB*, IMSDRUGNEWS, IMSPATENTS, IMSPRODUCT, IMSRESEARCH, IPA, MEDLINE, MRCK*, PATDPASPC,
       PROMT, PROUSDDR, PS, RTECS*, SYNTHLINE, TOXCENTER, USAN, USPAT2,
       USPATFULL
          (*File contains numerically searchable property data)
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Absolute stereochemistry. Rotation (-).

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

1137 REFERENCES IN FILE CA (1907 TO DATE)

20 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

1167 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> file merck COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 8.09 8.31

FULL ESTIMATED COST

FILE 'MRCK' ENTERED AT 22:50:05 ON 20 JUL 2010 COPYRIGHT (C) 2006, 2010 Merck Sharp & Dohme Corp., a subsidiary of Merck & Co., Whitehouse S

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=> s 11 L2 1 L1

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MERCK Number

(MNO): 1403918 (RN): **163222-33-1** CAS Registry No.

MERCK Index Name (MIN): Ezetimibe

CA Index Name (CN): (3R, 4S)-1-(4-Fluorophenyl)-3-[(3S)-3-(4-fluorophenyl)-3-

hydroxypropyl]-4-(4-hydroxyphenyl)-2-azetidinone

(CN): Sch-58235 Drug Code(s)

Trade Name(s) (CN): Ezetrol (Merck/Schering-Plough Pharmaceuticals (joint venture of Merck & Co., Inc. and Schering-Plough

Corp.); Merck/Schering-Plough); Zetia

(Merck/Schering-Plough Pharmaceuticals (joint venture

of Merck & Co., Inc. and Schering-Plough Corp.);

Merck/Schering-Plough)

File Segment. (FS): Active Monographs Molecular Form. (MF): C24 H21 F2 N O3

Wgt Composition (COMP): C 70.40%, H 5.17%, F 9.28%, N 3.42%, O 11.72%.

Molecular Weight (MW): 409.43

(RE): Cholesterol absorption inhibitor. Prepn: S. B. References Rosenblum et al., WO 9508532; eidem, US 5767115 (1995, 1998 both to Schering); idem et al., J. Med. Chem. 41, 973 (1998) DOI: 10.1021/jm970701f PMID: 9526571. Enantioselective synthesis: G. Wu et al., J. Org. Chem. 64, 3714 (1999) DOI: 10.1021/jo990428k PMID: 11674502. Activity in animals: M. van Heek et al., J. Pharmacol. Exp. Ther. 283, 157 (1997) PMID: 9336320. Metabolism and distribution: eldem, Br. J. Pharmacol. 129, 1748 (2000) DOI: 10.1038/sj.bjp.0703235 PMID: 10780982. Review of pharmacology and clinical studies: H. Bays, Expert Opin. Invest. Drugs 11, 1587-1604 (2002) DOI: 10.1517/13543784.11.11.1587 PMID: 12437505.

Absolute stereochemistry. Rotation (-).

Melting Point (MP):

Value MP deg C

======= 164 - 166

Optical Rotatory Power (ORP):

|Spectral| Value | Temp. | Line | ORP | ORP.T| ORP.SL | Note deg |deg C| | ____+__+___ -33.9|22|D|(c = 3 in methanol)

Other Properties (OCPP):

White solid, mp $164-166^{\circ}$. [α]D22 -33.9° (c = 3 in

methanol) .

Therapeutic Codes (THER):

Antilipemic.; Antilipemic; Others Referenced Patent (RPN):

WO9508532; US5767115

=> log y COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE TOTAL ENTRY SESSION 4.06 12.37

STN INTERNATIONAL LOGOFF AT 22:52:21 ON 20 JUL 2010